(FILE 'HOME' ENTERED AT 11:29:32 ON 12 AUG 2004)

	FILE	'REGI	STRY	ENT:	ERED	TA	11:32	2:05	ON	12	AUG	20	04					
L1			STRU	JCTUR:	E UPL	OAD	ED											
L2		15	S L	SSS	SAM													
L3		402	S L	SSS	FULL													
											3.00				017			0004
	FILE	' CAPL	US, N	JEDLTI	NE, U	SPA	TFULI	' E	NTER	RED	AT.	TT:	33:	23 (ON	12	AUG	2004
L4		6750	S L3	}														
L5		96	S L4	AND	(EYE	OR	GLAU	JCOM.	A OF	II S	VTRA	OCU	LAR	PR:	ESS	URE	3)	
L6		5	S L	AND	GLAU	COM	A											
L7		92	S L	AND	EYE													
L8		1	S L7	AND	INTR	AOC	ULAR	PRE	SSUR	₹E								

```
68
67
67
67
chain nodes :
                                      26
                                          27
                                              28
                                                  29
                                                     30
                                                                 33
                   17
                      18
                          21
                              22
                                  24
   12 13 14
              15
                              51
                                  58
                                      59
                                          60
                                              63
                                                  64
                                                      65
                                                         66
                                                             67
                                                                 68
                                                                    69
                                                                        70
                                                                            71
                                                                                72
                                                                                    78
   39 40
          41
              47
                  48
                      49
                          50
ring nodes :
                                     52
                                        53
                                             54
                                                55
                                                        57
   1 2 3 4
              5 6 7
                       8
                          9
                             10
                                 11
chain bonds :
                                     9-22 10-18 10-21 11-12 12-13 13-15 15-28 15-29
   1-8 2-14 4-24 5-26 6-27 9-17
                                     36-37
   15-47 30-31 30-32 33-34
                              33-35
                                           36-38 39-40 39-41 47-48 48-49 48-50 48-51
   50-57 58-59 59-60 63-64
                              65-66
                                     66-78
                                           67-68
                                                 67-69 70-71 70-72
```

```
ring nodes:

1 2 3 4 5 6 7 8 9 10 11 52 53 54 55 56 57

chain bonds:

1-8 2-14 4-24 5-26 6-27 9-17 9-22 10-18 10-21 11-12 12-13 13-15 15-28 15-29 15-47 30-31 30-32 33-34 33-35 36-37 36-38 39-40 39-41 47-48 48-49 48-50 48-51 50-57 58-59 59-60 63-64 65-66 66-78 67-68 67-69 70-71 70-72

ring bonds:

1-2 1-6 2-3 3-4 4-5 5-6 7-8 7-11 8-9 9-10 10-11 52-53 52-57 53-54 54-55 55-56 56-57

exact/norm bonds:

1-2 1-6 1-8 2-3 2-14 3-4 4-5 4-24 5-6 5-26 6-27 7-8 7-11 8-9 9-10 9-17 9-22 10-11 10-18 10-21 11-12 12-13 13-15 15-28 15-29 15-47 30-31 30-32 33-34 33-35 36-37 36-38 39-40 39-41 47-48 48-49 48-50 48-51 50-57 52-53 52-57 53-54 54-55 55-56 56-57 58-59 59-60 63-64 65-66 66-78 67-68 67-69 70-71 70-72
```

```
G1:H,F
```

G2:H,F,OH

G3:0,S

G4:H,Cl,Br,F,I,CN,Ak

G5:0,CH2,NH,[*1],[*2],[*3],[*4]

C:\Program Files\Stnexp\Queries\10087551-2.str

G6:CH2,[*5]

G7:H, [*6], [*7]

```
G8:CH2,[*1],[*4],[*8],[*9]
```

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom 11:Atom 12:CLASS 13:CLASS 14:CLASS 15:CLASS 17:CLASS 18:CLASS 21:CLASS 22:CLASS 24:CLASS 26:CLASS 27:CLASS 28:CLASS 29:CLASS 30:CLASS 31:CLASS 32:CLASS 33:CLASS 34:CLASS 35:CLASS 36:CLASS 37:CLASS 38:CLASS 39:CLASS 40:CLASS 41:CLASS 47:CLASS 48:CLASS 49:CLASS 50:CLASS 51:CLASS 52:Atom 53:Atom 54:Atom 55:Atom 56:CLASS 57:Atom 58:CLASS 59:CLASS 60:CLASS 63:CLASS 64:CLASS 65:CLASS 66:CLASS 67:CLASS 68:CLASS 69:CLASS 70:CLASS 71:CLASS 72:CLASS 78:CLASS

	(FILE 'HOME' ENTERED AT 09:00:47 ON 12 AUG 2004)
L1 L2 L3	FILE 'REGISTRY' ENTERED AT 09:01:04 ON 12 AUG 2004 STRUCTURE UPLOADED 15 S L1 SSS SAM 402 S L1 SSS FULL
	FILE 'CAPLUS, MEDLINE, USPATFULL' ENTERED AT 09:02:51 ON 12 AUG 2004
L4	6750 S L3
L5	96 S L4 AND (GLAUCOMA OR EYE OR INTRAOCULAR PRESSURE)
L6	5 S L5 AND GLAUCOMA

(FILE 'HOME' ENTERED AT 09:00:47 ON 12 AUG 2004)

	FILE 'REGISTRY' ENTERED AT 09:01:04 ON 12 AUG 2004
L1	STRUCTURE UPLOADED
L2	15 S L1 SSS SAM
L3	402 S L1 SSS FULL
	FILE 'CAPLUS, MEDLINE, USPATFULL' ENTERED AT 09:02:51 ON 12 AUG 2004
L4	6750 S L3
L5	96 S L4 AND (GLAUCOMA OR EYE OR INTRAOCULAR PRESSURE)
L6	5 S L5 AND GLAUCOMA
L7	3 S L5 AND HYPERTENSION

```
3 004
                             F C1 C1 F G 3 G 4
                                          6
                                          0 7 A
chain nodes :
   12 13 14 15 17
                      18
                          21
                              22
                                 24
                                     25
                                         27
                                             28
                                                 29
                                                     30
                                                         31
                                                             32
                                                                33
                                                                    34
          41
              42
                  48
                      49
                          50
                              51
                                 52
                                     59
                                         60
                                             61
                                                 64
                                                     65
                                                         66
                                                             67
                                                                 68
                                                                    69
                                                                        70
                                                                            71
   79
ring nodes :
   1 2 3 4
              5 6 7
                       8
                          9
                            10
                                11
                                     53
                                        54
                                            55
                                                56
                                                    57
                                                        58
chain bonds :
                                                        11-12 12-13 13-15
                                                                            15-29 15-30
   1-8 2-14 4-24 5-27 6-28 9-17
                                     9-22 10-18 10-21
   15-48 24-25 31-32 31-33 34-35
                                    34-36 37-38 37-39 40-41 40-42 48-49
                                                                            49-50 49-51
   49-52 51-58 59-60
                      60-61
                             64-65
                                     66-67 67-79 68-69 68-70 71-72
                                                                     71-73
```

G1:H,F G2:H,F,OH G3:O,S G4:H,Cl,Br,F,I,CN,Ak G5:O,CH2,NH,[*1],[*2],[*3],[*4] G6:CH2,[*5]

3-4 4-5 5-6

7-8

7-11 8-9

10-21 11-12 12-13 13-15 15-29 15-30 15-48

34-35 34-36 37-38 37-39 40-41 40-42 48-49 49-50 49-51 49-52 51-58 53-54 53-58 54-55 55-56 56-57 57-58 59-60 60-61 64-65 66-67 67-79 68-69 68-70 71-72 71-73

2-14 3-4 4-5 4-24 5-6 5-27

9-10 10-11 53-54 53-58

54-55

24-25 31-32 31-33

6-28 7-8 7-11 8-9 9-10 9-17

ring bonds :

G7:H,[*6],[*7]

1-2 1-6 2-3

1-2 1-6 1-8 2-3

9-22 10-11 10-18

56-57 57-58 exact/norm bonds :

C:\Program Files\Stnexp\Queries\10087551.str

G8:CH2,[*1],[*4],[*8],[*9]

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom 11:Atom 12:CLASS 13:CLASS 14:CLASS 15:CLASS 17:CLASS 18:CLASS 21:CLASS 22:CLASS 24:CLASS 25:CLASS 27:CLASS 28:CLASS 29:CLASS 30:CLASS 31:CLASS 32:CLASS 33:CLASS 34:CLASS 35:CLASS 36:CLASS 37:CLASS 38:CLASS 39:CLASS 40:CLASS 41:CLASS 42:CLASS 48:CLASS 49:CLASS 50:CLASS 51:CLASS 52:CLASS 53:Atom 54:Atom 55:Atom 56:Atom 57:CLASS 58:Atom 59:CLASS 70:CLASS 71:CLASS 72:CLASS 73:CLASS 79:CLASS 79:CLASS

```
ANSWER 1 OF 1 REGISTRY COPYRIGHT 2004 ACS on STN
L3
     133-89-1 REGISTRY
RN
     Uridine 5'-(trihydrogen diphosphate), P'-α-D-glucopyranosyl ester
           (CA INDEX NAME)
OTHER CA INDEX NAMES:
     Uridine 5'-(trihydrogen pyrophosphate), mono-\alpha-D-glucopyranosyl
     ester (8CI)
     Uridine 5'-pyrophosphate, \alpha-D-glucopyranosyl ester (6CI, 7CI)
OTHER NAMES:
     UDP-\alpha-D-Glucose
CN
CN
     UDP-D-glucose
CN
     UDP-Glc
CN
     UDP-Glucose
CN
     UDPG
CN
     Uridine 5'-(\alpha-D-glucopyranosyl pyrophosphate)
CN
     Uridine 5'-(trihydrogen pyrophosphate), mono-D-qlucosyl ester
CN
     Uridine 5'-diphosphate glucose
CN
     Uridine 5'-diphospho-α-D-glucose
     Uridine 5'-diphosphoglucose
CN
     Uridine diphosphate glucose
CN
     Uridine diphospho-D-glucose
CN
CN
     Uridine diphosphoglucose
     Uridine pyrophosphate-glucose
CN
FS
     STEREOSEARCH
     6659-38-7, 9002-11-3, 58-99-1, 73-37-0, 528-05-2, 25360-00-3, 99020-05-0,
DR
     99211-62-8, 30142-51-9, 30323-28-5
MF
     C15 H24 N2 O17 P2
CI
     COM
                  AGRICOLA, ANABSTR, BEILSTEIN*, BIOBUSINESS, BIOSIS,
LС
     STN Files:
       BIOTECHNO, CA, CANCERLIT, CAOLD, CAPLUS, CASREACT, CEN, CHEMINFORMRX.
       CHEMLIST, CSCHEM, DDFU, DRUGU, EMBASE, IFICDB, IFIPAT, IFIUDB, IPA,
       MEDLINE, MRCK*, NAPRALERT, PIRA, TOXCENTER, USPAT2, USPATFULL
         (*File contains numerically searchable property data)
     Other Sources:
                      EINECS**
         (**Enter CHEMLIST File for up-to-date regulatory information)
       CAplus document type: Conference; Dissertation; Journal; Patent; Report
DT.CA
       Roles from patents: ANST (Analytical study); BIOL (Biological study);
RL.P
       CMBI (Combinatorial study); FORM (Formation, nonpreparative); OCCU
       (Occurrence); PREP (Preparation); PROC (Process); PRP (Properties); RACT
       (Reactant or reagent); USES (Uses); NORL (No role in record)
       Roles for non-specific derivatives from patents: BIOL (Biological
RLD.P
       study)
      Roles from non-patents: ANST (Analytical study); BIOL (Biological
RL.NP
       study); FORM (Formation, nonpreparative); OCCU (Occurrence); PREP
       (Preparation); PROC (Process); PRP (Properties); RACT (Reactant or
       reagent); USES (Uses); NORL (No role in record)
RLD.NP Roles for non-specific derivatives from non-patents: BIOL (Biological
       study); FORM (Formation, nonpreparative); PREP (Preparation); PROC
       (Process); PRP (Properties); USES (Uses)
```

Absolute stereochemistry.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

- 2297 REFERENCES IN FILE CA (1907 TO DATE)
 - 23 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
- 2302 REFERENCES IN FILE CAPLUS (1907 TO DATE)
- 103 REFERENCES IN FILE CAOLD (PRIOR TO 1967)